FILE 'USPATFULL, CAPLUS, CAOLD' ENTERED AT 11:21:35 ON 12 JUN 2002

L1	545	S	95-05-6/RN OR 315-37-7/RN
L2	6	S	SULFIRAM OR (TETRAETHYLTHIODICARBONIC (W) DIAMIDE) OR ((BISDI
L3	141	S	MONOSULFIRAM OR SANIGAL OR CARCOCIDE OR SULFIRAMUM OR TETMOS
L4	691	S	(TESTOSTERONE (W) (ENANTHATE OR HEPTYLATE OR OENANTHATE OR HE
L5	993	S	L1 OR L2 OR L3 OR L4
L6	10783	S	3380-34-5/RN OR 1321-10-4/RN OR 59-50-7/RN OR 93-60-7/RN OR 4
L7	9617	S	TRICLOSAN OR CHLOROCRESOL OR CHLOROMETHYLPHENOL OR MONOCHLORO
L8	7600	S	ALIMEMAZINE OR ALIMEZINE OR METHYLPROMAZINE OR TERALEN OR OXY
L9	20274	S	L6 OR L7 OR L8
L10	21216	S	L9 OR L5
L11	64348	S	437-38-7/RN OR 27220-47-9/RN OR 65277-42-1/RN OR 12650-69-0/R
L12	195682	S	FENTANYL OR FENTANEST OR FENTANIL OR PHENTANYL OR ECONAZOLE O
L13	3560	S	TETRAMISOLE OR TETRAMISOL OR CHLORBUTANOL OR ACETOCHLORONE OR
L14	41513	S	TRICHLOROMETHYLPROPANOL OR TRICHLORODIMETHYLEHANOL OR (TRICHL
L15	4962	S	KETOPROFENE OR KETOPROPHEN OR FENOPROFEN OR ((PHENOXYPHENYLPR
L16	246802	S	L11 OR L12 OR L13 OR L14 OR L15
L17	264105	S	L9 OR L16
L18	63938	S	EUTECTIC
L19	366976	S	EMULSION

1,2-4

L2 ANSWER 1 OF 2 REGISTRY COPYRIGHT 2001 ACS
RN 1321-10-4 REGISTRY
CN Phenol, chloromethyl- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Cresol, chloro- (7CI, 8CI)
OTHER NAMES:
CN **Chlorocresol**
CN Chloromethylphenol
CN Monochlorocresol
DR 86006-41-9, 29468-35-7, 31308-59-5
MF C7 H7 Cl O
CI IDS, COM
LC STN Files: AGRICOLA, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CAOLD, CAPLUS, CHEMLIST, CIN, EMBASE, PROMT, TOXLINE, TOXLIT, USPATFULL
Other Sources: EINECS**
(**Enter CHEMLIST File for up-to-date regulatory information)

D1—Cl

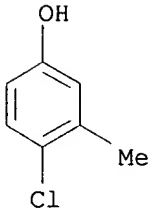
D1—OH

D1—Me

125 REFERENCES IN FILE CA (1967 TO DATE)
1 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
125 REFERENCES IN FILE CAPLUS (1967 TO DATE)
4 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

L2 ANSWER 2 OF 2 REGISTRY COPYRIGHT 2001 ACS
RN 59-50-7 REGISTRY
CN Phenol, 4-chloro-3-methyl- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN m-Cresol, 4-chloro- (8CI)
OTHER NAMES:
CN 1-Chloro-2-methyl-4-hydroxybenzene
CN 2-Chloro-5-hydroxytoluene
CN 3-Methyl-4-chlorophenol
CN 4-Chloro-3-cresol
CN 4-Chloro-3-methylphenol
CN 4-Chloro-5-methylphenol
CN 4-Chloro-m-cresol
CN 6-Chloro-3-hydroxytoluene
CN Aptal
CN Baktol
CN Baktolan
CN Candaseptic

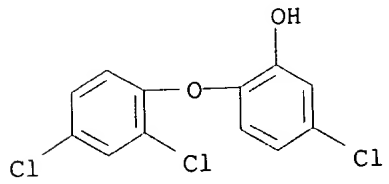
CN **Chlorocresol**
CN Neopredisan
CN Ottafact
CN p-Chloro-m-cresol
CN para-Chloro-meta-cresol
CN Parol
CN PCMC
CN Peritonan
CN Preventol CMK
CN Raschit
CN Raschit K
CN Rasen-Anicon
FS 3D CONCORD
MF C7 H7 Cl O
CI COM
LC STN Files: AGRICOLA, AIDSLINE, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DRUGU, EMBASE, GMELIN*, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IMSDIRECTORY, IPA, MEDLINE, MRCK*, MSDS-OHS, NIOSHTIC, PIRA, PROMT, RTECS*, SPECINFO, SYNTHLINE, TOXLINE, TOXLIT, ULIDAT, USAN, USPATFULL
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)



1268 REFERENCES IN FILE CA (1967 TO DATE)
14 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
1270 REFERENCES IN FILE CAPLUS (1967 TO DATE)
22 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

1,2-4

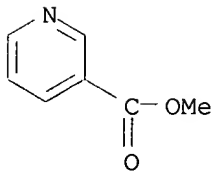
L1 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 3380-34-5 REGISTRY
CN Phenol, 5-chloro-2-(2,4-dichlorophenoxy)- (7CI, 8CI, 9CI) (CA INDEX NAME)
OTHER NAMES:
CN 2',4',4-Trichloro-2-hydroxydiphenyl ether
CN 2',4,4'-Trichloro-2-hydroxydiphenyl ether
CN 2,2'-Oxybis(1',5'-dichlorophenyl-5-chlorophenol)
CN 2,4,4'-Trichloro-2'-hydroxydiphenyl ether
CN 2-Hydroxy-2',4,4'-trichlorodiphenyl ether
CN 3-Chloro-6-(2,4-dichlorophenoxy)phenol
CN 4-Chloro-2-hydroxyphenyl 2,4-dichlorophenyl ether
CN 5-Chloro-2-(2,4-dichlorophenoxy)phenol
CN Bacti-Stat soap
CN CH 3565
CN DP 300
CN Irgacide LP 10
CN Irgasan
CN Irgasan CH 3565
CN Irgasan DP 30
CN Irgasan DP 300
CN Irgasan DP 3000
CN Irgasan PE 30
CN Irgasan PG 60
CN Microban Additive B
CN NM 100
CN THDP
CN **Triclosan**
CN Ultrafresh NM 100
CN Zilesan UW
FS 3D CONCORD
DR 112099-35-1, 88032-08-0
MF C12 H7 Cl3 O2
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DIOGENES, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS, NIOSHTIC, PIRA, PROMT, RTECS*, SPECINFO, TOXLINE, TOXLIT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)



1284 REFERENCES IN FILE CA (1967 TO DATE)
18 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
1286 REFERENCES IN FILE CAPLUS (1967 TO DATE)
2 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

1.2-4

L3 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 93-60-7 REGISTRY
CN 3-Pyridinecarboxylic acid, methyl ester (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Nicotinic acid, methyl ester (6CI, 7CI, 8CI)
OTHER NAMES:
CN 3-(Carbomethoxy)pyridine
CN 3-(Methoxycarbonyl)pyridine
CN m-(Methoxycarbonyl)pyridine
CN Methyl 3-pyridinecarboxylate
CN **Methyl nicotinate**
CN Nicometh
FS 3D CONCORD
DR 123574-61-8
MF C7 H7 N O2
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DIOGENES, DRUGU, EMBASE, HODOC*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS, NAPRALERT, NIOSHTIC, PROMT, RTECS*, SPECINFO, SYNTHLINE, TOXLINE, TOXLIT, USPATFULL
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)

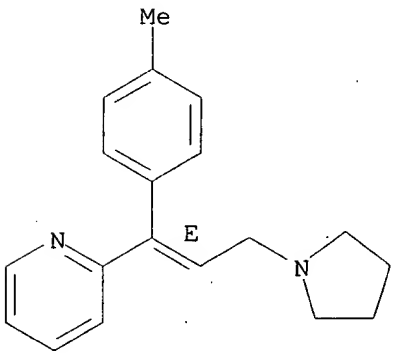


647 REFERENCES IN FILE CA (1967 TO DATE)
2 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
648 REFERENCES IN FILE CAPLUS (1967 TO DATE)
45 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

1, 2 = 4

L4 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 486-12-4 REGISTRY
CN Pyridine, 2-[(1E)-1-(4-methylphenyl)-3-(1-pyrrolidinyl)-1-propenyl]-
(9CI)
(CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Pyridine, 2-[1-(4-methylphenyl)-3-(1-pyrrolidinyl)-1-propenyl]-, (E)-
CN Pyridine, 2-[3-(1-pyrrolidinyl)-1-p-tolylpropenyl]-, (E)- (8CI)
OTHER NAMES:
CN trans-1-(2-Pyridyl)-3-pyrrolidino-1-p-tolylprop-1-ene
CN trans-1-(4-Methylphenyl)-1-(2-pyridyl)-3-pyrrolidinoprop-1-ene
CN trans-2-[3-(1-Pyrrolidinyl)-1-p-tolypropenyl]pyridine
CN Triprolidin
CN **Triprolidine**
CN Tripyrolidine
FS STEREOSEARCH
MF C19 H22 N2
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CHEMLIST, CIN, CSCHEM, DDFU,
DETHERM*, DIOGENES, DRUGU, EMBASE, HSDB*, IFICDB, IFIUDB, IPA, MEDLINE,
MRCK*, PROMT, RTECS*, SPECINFO, TOXLINE, TOXLIT, USAN, USPATFULL
(*File contains numerically searchable property data)
Other Sources: EINECS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)

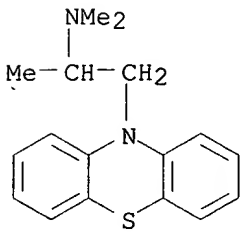
Double bond geometry as shown.



288 REFERENCES IN FILE CA (1967 TO DATE)
5 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
288 REFERENCES IN FILE CAPLUS (1967 TO DATE)
3 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

1, 2-4

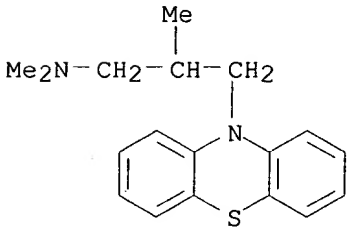
L5 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 60-87-7 REGISTRY
CN 10H-Phenothiazine-10-ethanamine, N,N,.alpha.-trimethyl- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Phenothiazine, 10-[2-(dimethylamino)propyl]- (8CI)
OTHER NAMES:
CN (.+-.)-Promethazine
CN (2-Dimethylamino-2-methyl)ethyl-N-dibenzoparathiazine
CN 10-[2-(Dimethylamino)propyl]phenothiazine
CN Dimapp
CN Diphergan
CN Hiberna
CN Proazamine
CN Procit
CN **Promethazine**
CN Protazine
CN Prothazin
CN RP 3277
CN Vallergine
FS 3D CONCORD
DR 73745-50-3
MF C17 H20 N2 S
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU, DIOGENES, DRUGU, EMBASE, GMELIN*, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IMSDIRECTORY, IPA, MEDLINE, MRCK*, NIOSHTIC, PROMT, RTECS*, SPECINFO, TOXLINE, TOXLIT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: EINECS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)



2043 REFERENCES IN FILE CA (1967 TO DATE)
42 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
2047 REFERENCES IN FILE CAPLUS (1967 TO DATE)
43 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

1, 2 - 4

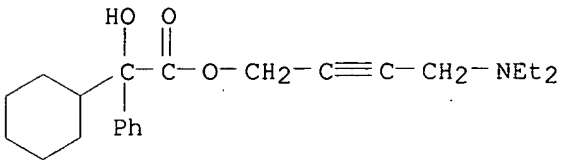
L6 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 84-96-8 REGISTRY
CN 10H-Phenothiazine-10-propanamine, N,N,.beta.-trimethyl- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Phenothiazine, 10-[3-(dimethylamino)-2-methylpropyl]- (6CI, 8CI)
OTHER NAMES:
CN (.+-.)-Alimemazine
CN (.+-.)-Trimeprazine
CN 10-(2-Methyl-3-dimethylaminopropyl)phenothiazine
CN 10-[3-(Dimethylamino)-2-methylpropyl]phenothiazine
CN Alimemazine
CN Alimezine
CN Bayer 1219
CN dl-Trimeprazine
CN Methylpromazine
CN Teralen
CN **Trimeprazine**
FS 3D CONCORD
DR 35309-60-5, 47138-21-6
MF C18 H22 N2 S
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CHEMLIST, CIN, DDFU, DRUGU, EMBASE, HSDB*, IPA, MEDLINE, MRCK*, NIOSHTIC, RTECS*, SPECINFO, TOXLINE, TOXLIT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: EINECS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)



406 REFERENCES IN FILE CA (1967 TO DATE)
7 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
407 REFERENCES IN FILE CAPLUS (1967 TO DATE)
48 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

1,2-4

L8 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 5633-20-5 REGISTRY
CN Benzeneacetic acid, .alpha.-cyclohexyl-.alpha.-hydroxy-, 4-(diethylamino)-2-butynyl ester (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN 2-Butyn-1-ol, 4-(diethylamino)-, .alpha.-phenylcyclohexaneglycolate (ester)
CN Cyclohexaneglycolic acid, .alpha.-phenyl-, 4-(diethylamino)-2-butynyl ester (8CI)
OTHER NAMES:
CN (.+-.)-Oxybutynin
CN (RS)-Oxybutynin
CN 4-Diethylamino-2-butynyl .alpha.-phenylcyclohexaneglycolate
CN **Oxybutynin**
FS 3D CONCORD
DR 119579-36-1
MF C22 H31 N O3
CI COM
LC STN Files: ADISINSIGHT, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CANCERLIT, CAPLUS, CBNB, CEN, CHEMCATS, CIN, CSCHEM, DDFU, DIOGENES, DRUGU, EMBASE, HSDB*, IPA, MEDLINE, MRCK*, PHAR, PROMT, SPECINFO, TOXLINE, TOXLIT, USAN, USPATFULL
(*File contains numerically searchable property data)
Other Sources: WHO

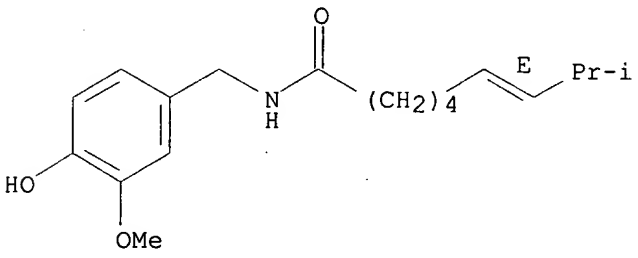


181 REFERENCES IN FILE CA (1967 TO DATE)
3 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
183 REFERENCES IN FILE CAPLUS (1967 TO DATE)

1,2-4

L9 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 404-86-4 REGISTRY
CN 6-Nonenamide, N-[(4-hydroxy-3-methoxyphenyl)methyl]-8-methyl-, (6E)-(9CI)
(CA INDEX NAME)
OTHER CA INDEX NAMES:
CN 6-Nonenamide, 8-methyl-N-vanillyl-, (E)- (8CI)
CN 6-Nonenamide, N-[(4-hydroxy-3-methoxyphenyl)methyl]-8-methyl-, (E)-
CN **Capsaicin (6CI)**
OTHER NAMES:
CN (E)-N-(4-Hydroxy-3-methoxybenzyl)-8-methylnon-6-enamide
CN Capsaicine
CN Ratden PE 40
CN Zostrix
FS STEREOSEARCH
MF C18 H27 N O3
CI COM
LC STN Files: ADISINSIGHT, AGRICOLA, AIDSLINE, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DRUGNL, DRUGU, DRUGUPDATES, EMBASE, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS, NAPRALERT, NIOSHTIC, PHAR, PIRA, PROMT, RTECS*, TOXLINE, TOXLIT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)

Double bond geometry as shown.



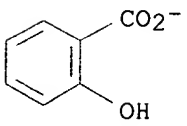
2880 REFERENCES IN FILE CA (1967 TO DATE)
58 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

1, 2 - 4

L11 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 2016-36-6 REGISTRY
CN Ethanaminium, 2-hydroxy-N,N,N-trimethyl-, salt with 2-hydroxybenzoic acid (1:1) (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Benzoic acid, 2-hydroxy-, ion(1-), 2-hydroxy-N,N,N-trimethylethanaminium (9CI)
CN **Choline salicylate (6CI)**
CN Choline, salicylate (salt) (7CI, 8CI)
CN Salicylic acid, ion(1-), choline (8CI)
OTHER NAMES:
CN (2-Hydroxyethyl)trimethylammonium salicylate
CN Actasal
CN Arret
CN Arthropan
CN Artrobione
CN Mundisal
CN Salicol
CN Satibon
CN Syrap
DR 54391-51-4
MF C7 H5 O3 . C5 H14 N O
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CBNB, CHEMLIST, CIN, CSCHEM, DDFU, DIOGENES, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IMSDIRECTORY, IPA, MEDLINE, MRCK*, PROMT, RTECS*, TOXLINE, TOXLIT, USAN, USPATFULL
(*File contains numerically searchable property data)
Other Sources: EINECS**, NDSL**, TSCA**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)

CM 1

CRN 63-36-5
CMF C7 H5 O3



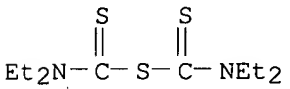
CM 2

CRN 62-49-7
CMF C5 H14 N O

$Me_3^+N-CH_2-CH_2-OH$

67 REFERENCES IN FILE CA (1967 TO DATE)
4 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
67 REFERENCES IN FILE CAPLUS (1967 TO DATE)
14 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

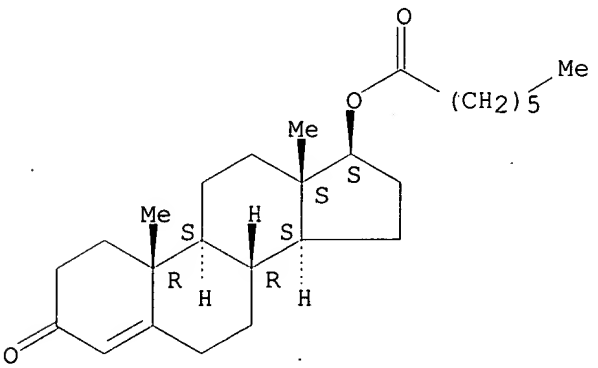
L7 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 95-05-6 REGISTRY
CN Thiodicarbonic diamide ([(H2N)C(S)]2S), tetraethyl- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Sulfide, bis(diethylthiocarbamoyl) (6CI, 7CI, 8CI)
OTHER NAMES:
CN Bis(diethylthiocarbamoyl) sulfide
CN Bis(diethylthiocarbamyl) sulfide
CN Bis(N,N-diethylthiocarbamoyl) sulfide
CN Carbamodithioic acid, diethyl-, anhydrosulfide
CN Kutka
CN Methanethioamide, 1,1'-thiobis[N,N-diethyl-
CN Monosulfiram
CN Sanigal
CN Sarcocide B
CN Sulfide, bis[(diethylamino)thioxomethyl]
CN **Sulfiram**
CN Sulfirame
CN Sulfiramum
CN Tetmos
CN Tetmosol
CN Tetraethylthiuram monosulfide
CN Tetrucid
FS 3D CONCORD
MF C10 H20 N2 S3
CI COM
LC STN Files: AGRICOLA, BEILSTEIN*, BIOSIS, CA, CABA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMLIST, DDFU, DRUGU, EMBASE, HODOC*, IMSDIRECTORY, IPA, MEDLINE, MRCK*, PROMT, RTECS*, TOXLINE, TOXLIT, USAN, USPATFULL
(*File contains numerically searchable property data)
Other Sources: EINECS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)



37 REFERENCES IN FILE CA (1967 TO DATE)
37 REFERENCES IN FILE CAPLUS (1967 TO DATE)
8 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

L10 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 315-37-7 REGISTRY
CN Androst-4-en-3-one, 17-[(1-oxoheptyl)oxy]-, (17.beta.)- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Testosterone, heptanoate (6CI, 8CI)
OTHER NAMES:
CN 17.beta.-Enanthoxyandrost-4-en-3-one
CN 17.beta.-Hydroxyandrost-4-en-3-one enanthate
CN 4-Androsten-3-one 17.beta.-enanthate
CN Androtardyl
CN Delatestryl
CN Reposo TMD
CN Testenate
CN Testosterone 17-enanthate
CN **Testosterone enanthate**
CN Testosterone heptylate
CN Testosterone oenanthate
FS STEREOSEARCH
DR 11111-10-7
MF C26 H40 O3
CI COM
LC STN Files: ADISINSIGHT, AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS, CBNB, CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU, DIOGENES, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS, PROMT, RTECS*, TOXLINE, TOXLIT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: EINECS**
(**Enter CHEMLIST File for up-to-date regulatory information)

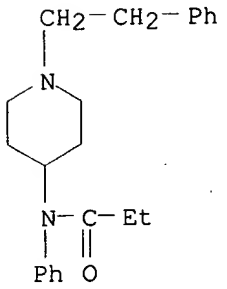
Absolute stereochemistry. Rotation (+).



389 REFERENCES IN FILE CA (1967 TO DATE)
389 REFERENCES IN FILE CAPLUS (1967 TO DATE)
45 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

2-4

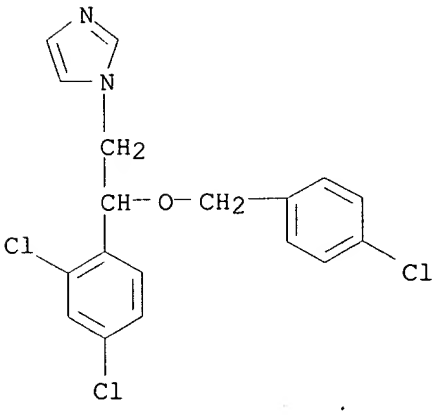
L17 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 437-38-7 REGISTRY
CN Propanamide, N-phenyl-N-[1-(2-phenylethyl)-4-piperidinyl]- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Propionanilide, N-(1-phenethyl-4-piperidyl)- (7CI, 8CI)
OTHER NAMES:
CN 1-Phenethyl-4-(N-phenylpropionamido)piperidine
CN 1-Phenethyl-4-N-propionylanilinopiperidine
CN Durogesic
CN Fentanest
CN Fentanil
CN **Fentanyl**
CN N-[1-(2-Phenylethyl)-4-piperidinyl]propionanilide
CN Phentanyl
CN R 4263
FS 3D CONCORD
DR 80832-90-2
MF C22 H28 N2 O
CI COM
LC STN Files: ADISINSIGHT, AGRICOLA, AIDSLINE, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS, CBNB, CEN, CHEMCATS, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DIOGENES, DRUGNL, DRUGU, DRUGUPDATES, EMBASE, HODOC*, HSDB*, IFICDB, IFIUDB, IMSDIRECTORY, IPA, MEDLINE, MRCK*, MSDS-OHS, NIOSHTIC, PHAR, PROMT, RTECS*, SPECINFO, TOXLINE, TOXLIT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: EINECS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)



2314 REFERENCES IN FILE CA (1967 TO DATE)
73 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
2315 REFERENCES IN FILE CAPLUS (1967 TO DATE)
20 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

2-4

L18 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 27220-47-9 REGISTRY
CN 1H-Imidazole,
1-[2-[(4-chlorophenyl)methoxy]-2-(2,4-dichlorophenyl)ethyl]-
(9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Imidazole, 1-[2,4-dichloro-.beta.-[(p-chlorobenzyl)oxy]phenethyl]- (8CI)
OTHER NAMES:
CN (.+-.)-Econazole
CN 1-[2,4-Dichloro-.beta.-[(p-chlorobenzyl)oxy]phenethyl]imidazole
CN **Econazole**
CN Spectazole
FS 3D CONCORD
DR 68797-30-8
MF C18 H15 Cl3 N2 O
CI COM
LC STN Files: ADISINSIGHT, AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CABA, CANCERLIT, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DIOGENES, DRUGPAT, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IMSDIRECTORY, IPA, MEDLINE, MRCK*, NIOSHTIC, PHAR, PROMT, RTECS*, TOXLINE, TOXLIT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: EINECS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)

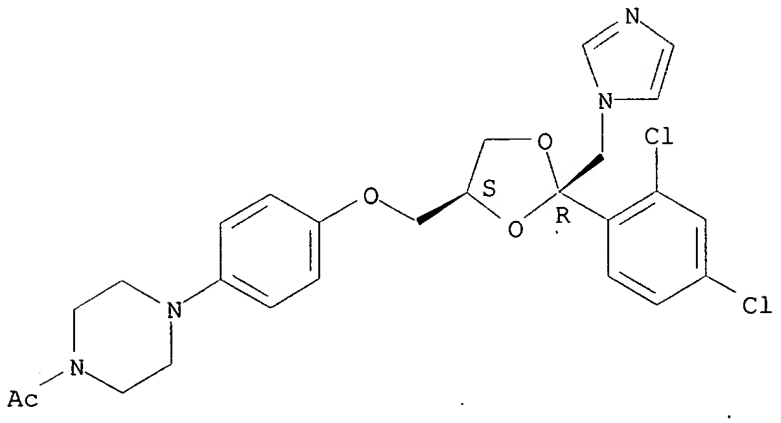


456 REFERENCES IN FILE CA (1967 TO DATE)
10 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
456 REFERENCES IN FILE CAPLUS (1967 TO DATE)

3-4

L19 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 65277-42-1 REGISTRY
CN Piperazine,
1-acetyl-4-[4-[[(2R,4S)-2-(2,4-dichlorophenyl)-2-(1H-imidazol-1-ylmethyl)-1,3-dioxolan-4-yl]methoxy]phenyl]-, rel- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Piperazine, 1-acetyl-4-[4-[[2-(2,4-dichlorophenyl)-2-(1H-imidazol-1-ylmethyl)-1,3-dioxolan-4-yl]methoxy]phenyl]-, cis-
OTHER NAMES:
CN (.+-.)-Ketoconazole
CN Fungoral
CN **Ketoconazole**
CN Nizoral
CN R 41400
FS STEREOSEARCH
DR 72093-26-6
MF C26 H28 Cl2 N4 O4
CI COM
LC STN Files: AGRICOLA, AIDSLINE, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
BIOTECHNO, CA, CABA, CANCERLIT, CAPLUS, CBNB, CEN, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DIOGENES, DRUGPAT, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IMSDIRECTORY, IPA, MEDLINE, MRCK*, MSDS-OHS, NIOSHTIC, PHAR, PROMT, RTECS*, SPECINFO, SYNTHLINE, TOXLINE, TOXLIT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: EINECS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)

Relative stereochemistry.

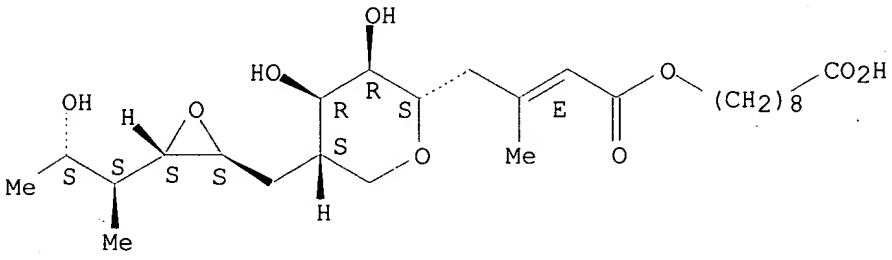


1863 REFERENCES IN FILE CA (1967 TO DATE)
33 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
1868 REFERENCES IN FILE CAPLUS (1967 TO DATE)

2-4

L20 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 12650-69-0 REGISTRY
CN L-talo-Non-2-enonic acid, 5,9-anhydro-2,3,4,8-tetradeoxy-8-[[(2S,3S)-3-[(1S,2S)-2-hydroxy-1-methylpropyl]oxiranyl]methyl]-3-methyl-, 8-carboxyoctyl ester, (2E)- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN L-talo-Non-2-enonic acid, 5,9-anhydro-2,3,4,8-tetradeoxy-8-[[3-(2-hydroxy-1-methylpropyl)oxiranyl]methyl]-3-methyl-, 8-carboxyoctyl ester, [2E,8[2S,3S(1S,2S)]]-
OTHER NAMES:
CN Bactroban
CN Bactroban Ointment
CN **Mupirocin**
CN Pseudomonic acid
CN Pseudomonic acid A
CN trans-Pseudomonic acid
FS STEREOSEARCH
DR 62916-63-6
MF C26 H44 O9
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CABA, CANCERLIT, CAPLUS, CBNB, CEN, CIN, DDFU, DIOGENES, DRUGNL, DRUGPAT, DRUGU, DRUGUPDATES, EMBASE, IFICDB, IFIPAT, IFIUDB, IMSDIRECTORY, IPA, MEDLINE, MRCK*, NAPRALERT, PHAR, PROMT, RTECS*, TOXLINE, TOXLIT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: WHO

Absolute stereochemistry.
Double bond geometry as shown.

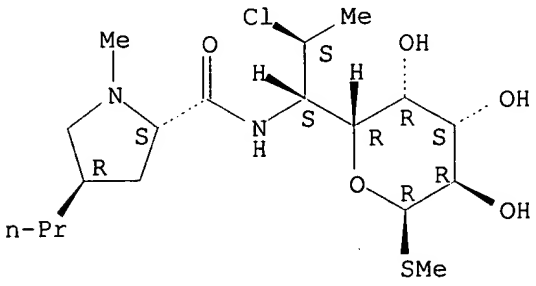


308 REFERENCES IN FILE CA (1967 TO DATE)
18 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
309 REFERENCES IN FILE CAPLUS (1967 TO DATE)

2 - 4

L21 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 18323-44-9 REGISTRY
CN L-threo-.alpha.-D-galacto-Octopyranoside, methyl 7-chloro-6,7,8-trideoxy-6-[[[(2S,4R)-1-methyl-4-propyl-2-pyrrolidinyl]carbonyl]amino]-1-thio- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN L-threo-.alpha.-D-galacto-Octopyranoside, methyl 7-chloro-6,7,8-trideoxy-6-[[(1-methyl-4-propyl-2-pyrrolidinyl)carbonyl]amino]-1-thio-, (2S-trans)-
CN L-threo-D-galacto-Octopyranoside, methyl 7-chloro-6,7,8-trideoxy-6-(1-methyl-4-propyl-L-2-pyrrolidinecarboxamido)-1-thio-, trans- .alpha.- (8CI)
OTHER NAMES:
CN 7(S)-Chloro-7-deoxylincomycin
CN 7-CDL
CN 7-Chloro-7-deoxylincomycin
CN 7-Chlorolincomycin
CN 7-Deoxy-7(S)-chlorolincomycin
CN Chlolincocin
CN Cleocin
CN **Clindamycin**
CN Clinimycin
CN Dalacin C
CN Sobelin
CN U 21251
CN U-21,251
FS STEREOSEARCH
DR 13441-63-9, 24620-78-8, 24696-19-3, 16669-21-9
MF C18 H33 Cl N2 O5 S
CI COM
LC STN Files: AGRICOLA, AIDSLINE, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
BIOTECHNO, CA, CABA, CANCERLIT, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU, DIOGENES, DRUGPAT, DRUGU, EMBASE, HSDB*, IFICDB, IFIPAT, IFIUDB, IMSDIRECTORY, IPA, MEDLINE, MRCK*, NAPRALERT, NIOSHTIC, PHAR, PROMT, RTECS*, TOXLINE, TOXLIT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: EINECS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)

Absolute stereochemistry.



2425 REFERENCES IN FILE CA (1967 TO DATE)
21 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
2429 REFERENCES IN FILE CAPLUS (1967 TO DATE)

2-4

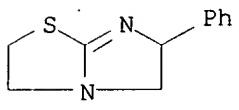
L22 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 7553-56-2 REGISTRY
CN **Iodine (8CI, 9CI)** (CA INDEX NAME)
OTHER NAMES:
CN Actomar
CN Diatomic iodine
CN Diiodine
CN Eranol
CN Iodel FD
CN Iodine (127I2)
CN Iodine colloidal
CN Iodine crystals
CN Iodine molecule (I2)
CN Iodine sublimed
CN Iosan Superdip
CN Jodosan
CN Molecular iodine
FS 3D CONCORD
DR 8012-81-5, 8012-85-9, 8031-47-8, 24503-90-0
MF I2
CI COM
LC STN Files: ADISINSIGHT, AGRICOLA, AIDSLINE, ANABSTR, APILIT, APILIT2, APIPAT, APIPAT2, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CABA, CANCERLIT, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CHEMINFORMRX, CHEMLIST, CHEMSAFE, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DIOGENES, DIPPR*, DRUGU, EMBASE, GMELIN*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS, NAPRALERT, NIOSHTIC, PDLCOM*, PHAR, PIRA, PROMT, RTECS*, TOXLINE, TOXLIT, TRCTHERMO*, TULSA, ULIDAT, USAN, USPATFULL, VETU, VTB
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)

I—I

35260 REFERENCES IN FILE CA (1967 TO DATE)
2274 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
35276 REFERENCES IN FILE CAPLUS (1967 TO DATE)

2 - 4/

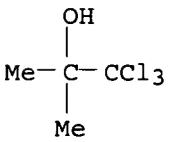
L23 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 5036-02-2 REGISTRY
CN Imidazo[2,1-b]thiazole, 2,3,5,6-tetrahydro-6-phenyl- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Imidazo[2,1-b]thiazole, 2,3,5,6-tetrahydro-6-phenyl-, (.+-.)- (8CI)
OTHER NAMES:
CN (.+-.)-2,3,5,6-Tetrahydro-6-phenylimidazo[2,1-b]thiazole
CN (.+-.)-Tetramisole
CN 2,3,5,6-Tetrahydro-6-phenylimidazo[2,1-b]thiazole
CN 6-Phenyl-2,3,5,6-tetrahydroimidazo[2,1-b]thiazole
CN dl-2,3,5,6-Tetrahydro-6-phenylimidazo(2,1-b)thiazole
CN dl-Tetramisol
CN dl-Tetramisole
CN Nilverm base
CN Tetramisol
CN **Tetramisole**
FS 3D CONCORD
DR 6649-23-6
MF C11 H12 N2 S
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, NIOSHTIC, PROMT, RTECS*, SPECINFO, TOXLINE, TOXLIT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: EINECS**, NDSL**, TSCA**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)



231 REFERENCES IN FILE CA (1967 TO DATE)
10 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
231 REFERENCES IN FILE CAPLUS (1967 TO DATE)
1 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

2 – 4/

L2 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 57-15-8 REGISTRY
CN 2-Propanol, 1,1,1-trichloro-2-methyl- (6CI, 8CI, 9CI) (CA INDEX NAME)
OTHER NAMES:
CN .beta.,.beta.,.beta.-Trichloro-tert-butyl alcohol
CN 1,1,1-Trichloro-2-methyl-2-propanol
CN 1,1,1-Trichloro-tert-butyl alcohol
CN 2,2,2-Trichloro-1,1-dimethylethanol
CN 2-(Trichloromethyl)-2-propanol
CN Acetochlorone
CN Acetonchloroform
CN Acetone chloroform
CN Anhydrous chlorobutanol
CN **Chlorbutanol**
CN Chlorbutol
CN Chloreton
CN Chloretone
CN Chlorobutanol
CN Chlortran
CN Clortran
CN Dentalone
CN Khloreton
CN Methaform
CN Sedaform
CN Trichloro-tert-butyl alcohol
FS 3D CONCORD
MF C4 H7 Cl3 O
CI COM
LC STN Files: ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DIOGENES, DRUGU, EMBASE, GMELIN*, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS, NAPRALERT, NIOSHTIC, PIRA, PROMT, RTECS*, SPECINFO, SYNTHLINE, TOXLIT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)

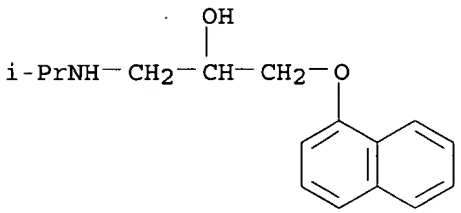


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

472 REFERENCES IN FILE CA (1967 TO DATE)
4 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
473 REFERENCES IN FILE CAPLUS (1967 TO DATE)
30 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

3 - 4

L3 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 525-66-6 REGISTRY
CN 2-Propanol, 1-[(1-methylethyl)amino]-3-(1-naphthalenyloxy)- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN 2-Propanol, 1-(isopropylamino)-3-(1-naphthyloxy)- (7CI, 8CI)
OTHER NAMES:
CN (.+-.)-Propranolol
CN .beta.-Propranolol
CN 1-(1-Naphthyloxy)-3-(isopropylamino)-2-propanol
CN 1-(Isopropylamino)-3-(1-naphthyloxy)-2-propanol
CN AY 64043
CN Betalong
CN dl-Propranolol
CN DL-Propranolol
CN **Propranolol**
CN Proprasylyt
CN Racemic propranolol
CN Reducor
FS 3D CONCORD
DR 13013-17-7
MF C16 H21 N O2
CI COM
LC STN Files: ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSNB, DDFU, DIOGENES, DRUGPAT, DRUGU, EMBASE, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, NIOSHTIC, PHAR, PHARMASEARCH, PROMT, RTECS*, SPECINFO, TOXLIT, ULIDAT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: EINECS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)

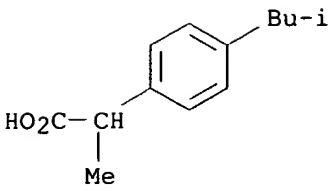


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

8877 REFERENCES IN FILE CA (1967 TO DATE)
102 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
8882 REFERENCES IN FILE CAPLUS (1967 TO DATE)
1 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

2 -4

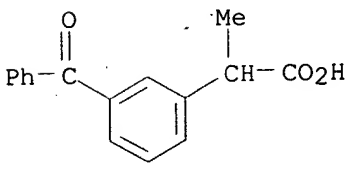
L12 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 15687-27-1 REGISTRY
CN Benzeneacetic acid, .alpha.-methyl-4-(2-methylpropyl)- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Hydratropic acid, p-isobutyl- (7CI, 8CI)
OTHER NAMES:
CN (.+-.)-.alpha.-Methyl-4-(2-methylpropyl)benzeneacetic acid
CN (.+-.)-2-(p-Isobutylphenyl)propionic acid
CN (.+-.)-Ibuprofen
CN (.+-.)-Ibuprophen
CN (4-Isobutylphenyl)-.alpha.-methylacetic acid
CN (RS)-Ibuprofen
CN (S)-4-Isobutyl-.alpha.-methylphenylacetic acid
CN .alpha.-(4-Isobutylphenyl)propionic acid
CN .alpha.-Methyl-4-(2-methylpropyl)benzeneacetic acid
CN 2-(4'-Isobutylphenyl)propionic acid
CN 2-(4-Isobutylphenyl)propanoic acid
CN 2-(p-Isobutylphenyl)propionic acid
CN 4-Isobutyl-.alpha.-methylphenylacetic acid
CN 4-Isobutylhydratropic acid
CN Advil
CN Brufen
CN dl-Ibuprofen
CN Ibufen
CN **Ibuprofen**
CN IP 82
CN Motrin
CN Nuprin
CN Nurofen
CN p-Isobutyl-2-phenylpropionic acid
CN p-Isobutylhydratropic acid
CN Paduden
CN Proflex
CN RD 13621
CN Rufin
CN Unipron
FS 3D CONCORD
DR 58560-75-1
MF C13 H18 O2
CI COM
LC STN Files: ADISINSIGHT, AGRICOLA, AIDSLINE, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, DDFU, DETHERM*, DIOGENES, DIPPR*, DRUGPAT, DRUGU, EMBASE, HSDB*, IFICDB, IFIPAT, IFIUDB, IMSDIRECTORY, IPA, MEDLINE, MRCK*, MSDS-OHS, NIOSHTIC, PHAR, PIRA, PROMT, RTECS*, SPECINFO, TOXLINE, TOXLIT, ULIDAT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)



5016 REFERENCES IN FILE CA (1967 TO DATE)
148 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
5023 REFERENCES IN FILE CAPLUS (1967 TO DATE)
2 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

2 ~4

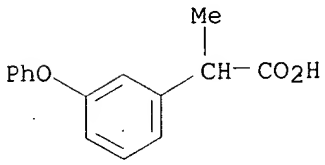
L13 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 22071-15-4 REGISTRY
CN Benzeneacetic acid, 3-benzoyl-.alpha.-methyl- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Hydratropic acid, m-benzoyl- (8CI)
OTHER NAMES:
CN (.+-.)-2-(3-Benzoylphenyl)propionic acid
CN (.+-.)-3-Benzoyl-.alpha.-methylbenzeneacetic acid
CN (.+-.)-Ketoprofen
CN (.+-.)-m-Benzoylhydratropic acid
CN (RS)-Ketoprofen
CN .alpha.-(3-Benzoylphenyl)propionic acid
CN 19583RP
CN 2-(3-Benzoylphenyl)propionic acid
CN 2-(m-Benzoylphenyl)propionic acid
CN 3-Benzoyl-.alpha.-methylbenzeneacetic acid
CN 3-Benzoylhydratropic acid
CN Alrheumun
CN Aneol
CN Capisten
CN Epatec
CN **Ketoprofen**
CN Ketoprofene
CN Ketoprophen
CN m-Benzoylhydratropic acid
CN Orudis
CN Oruvail
CN Profenid
CN R.P. 19583
CN Racemic ketoprofen
CN RU 4733
FS 3D CONCORD
DR 172964-50-0, 22161-86-0
MF C16 H14 O3
CI COM
LC STN Files: AGRICOLA, AIDSLINE, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
BIOTECHNO, CA, CABA, CANCERLIT, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DIOGENES, DRUGPAT, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IMSDIRECTORY, IPA, MEDLINE, MRCK*, MSDS-OHS, NIOSHTIC, PHAR, PROMT, RTECS*, SPECINFO, SYNTHLINE, TOXLINE, TOXLIT, ULIDAT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: EINECS**, NDSL**, TSCA**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)



2400 REFERENCES IN FILE CA (1967 TO DATE)
82 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
2407 REFERENCES IN FILE CAPLUS (1967 TO DATE)

2-4

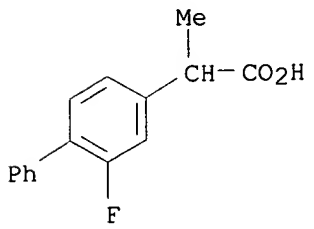
L14 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 29679-58-1 REGISTRY
CN Benzeneacetic acid, .alpha.-methyl-3-phenoxy- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Hydratropic acid, m-phenoxy- (8CI)
OTHER NAMES:
CN (.+-.)-2-(3-Phenoxyphenyl)propionic acid
CN (.+-.)-Fenoprofen
CN (.+-.)-m-Phenoxyhydratropic acid
CN .alpha.-Methyl-3-phenoxybenzeneacetic acid
CN 2-(3-Phenoxyphenyl)propionic acid
CN 2-(m-Phenoxyphenyl)propionic acid
CN 3-Phenoxyhydratropic acid
CN dl-2-(3-Phenoxyphenyl)propionic acid
CN **Fenoprofen**
FS 3D CONCORD
DR 31879-05-7
MF C15 H14 O3
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CANCERLIT, CAPLUS, CASREACT, CHEMCATS, CHEMLIST, CIN, DDFU, DRUGPAT, DRUGU, EMBASE, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, NIOSHTIC, PROMT, RTECS*, SPECINFO, TOXLINE, TOXLIT, ULIDAT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: EINECS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)



267 REFERENCES IN FILE CA (1967 TO DATE)
12 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
268 REFERENCES IN FILE CAPLUS (1967 TO DATE)

2 - 4

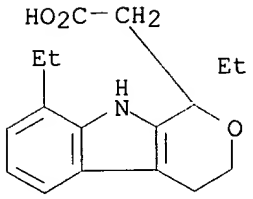
L15 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 5104-49-4 REGISTRY
CN [1,1'-Biphenyl]-4-acetic acid, 2-fluoro-.alpha.-methyl- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN 4-Biphenylacetic acid, 2-fluoro-.alpha.-methyl- (8CI)
CN Hydratropic acid, 3-fluoro-4-phenyl- (7CI)
OTHER NAMES:
CN (.+-.)-Flurbiprofen
CN 2-(2-Fluoro-4-biphenyl)propionic acid
CN 2-(2-Fluoro-4-biphenylyl)propanoic acid
CN 2-(2-Fluoro-4-biphenylyl)propionic acid
CN 2-Fluoro-.alpha.-methyl-4-biphenylacetic acid
CN 2-Fluoro-.alpha.-methyl-4-diphenylacetic acid
CN 3-Fluoro-4-phenylhydratropic acid
CN Ansaid
CN dl-2-(2-Fluoro-4-biphenylyl)propionic acid
CN dl-Flurbiprofen
CN Flugalin
CN **Flurbiprofen**
CN FP 70
CN FP-A
CN Froben
CN rac-Flurbiprofen
CN Racemic flurbiprofen
CN U 27182
FS 3D CONCORD
DR 51543-38-5, 79212-68-3
MF C15 H13 F O2
CI COM
LC STN Files: ADISINSIGHT, AGRICOLA, AIDSLINE, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, DDFU, DETHERM*, DIOGENES, DRUGPAT, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IMSDIRECTORY, IPA, MEDLINE, MRCK*, PHAR, PROMT, RTECS*, SPECINFO, SYNTHLINE, TOXLINE, TOXLIT, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)



1574 REFERENCES IN FILE CA (1967 TO DATE)
56 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
1577 REFERENCES IN FILE CAPLUS (1967 TO DATE)
1 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

2-4

L16 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2001 ACS
RN 41340-25-4 REGISTRY
CN Pyrano[3,4-b]indole-1-acetic acid, 1,8-diethyl-1,3,4,9-tetrahydro- (9CI) (CA INDEX NAME)
OTHER NAMES:
CN (.+-.)-Etodolac
CN (RS)-Etodolic acid
CN AY 24236
CN **Etodolac**
CN Etodolic acid
CN NIH 9918
FS 3D CONCORD
DR 87226-38-8
MF C17 H21 N O3
CI COM
LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CANCERLIT, CAPLUS, CASREACT, CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU, DIOGENES, DRUGNL, DRUGPAT, DRUGU, DRUGUPDATES, EMBASE, IFICDB, IFIPAT, IFIUDB, IMSDIRECTORY, IPA, MEDLINE, MRCK*, MSDS-OHS, PHAR, PROMT, RTECS*, SYNTHLINE, TOXLINE, TOXLIT, USAN, USPATFULL
(*File contains numerically searchable property data)
Other Sources: DSL**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)



351 REFERENCES IN FILE CA (1967 TO DATE)
25 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
352 REFERENCES IN FILE CAPLUS (1967 TO DATE)

L6 ANSWER 4 OF 6 USPATFULL

ACCESSION NUMBER: 93:104689 USPATFULL

TITLE: Sterols, their fatty acid esters and glucosides; processes for their preparation; spontaneously dispersible agents containing these compounds, and their use for treatment of tumors

INVENTOR(S): Eugster, Carl, Riehen, Switzerland
Eugster, Conrad, Wallisellen, Switzerland
Haldemann, Walter, Binningen, Switzerland
Rivara, Giorgio, Turin, Italy

PATENT ASSIGNEE(S): Marigen S.A., Riehen, Switzerland (non-U.S. corporation)

	NUMBER	DATE	
PATENT INFORMATION:	US 5270041	19931214	
	WO 9101139	19910207	
APPLICATION INFO.:	US 1991-634215	19910215	(7)
	WO 1990-CH164	19900706	
		19910215	PCT 371 date
		19910215	PCT 102(e) date

	NUMBER	DATE
PRIORITY INFORMATION:	CH 1989-2727	19890721
	CH 1989-4308	19891202

DOCUMENT TYPE: Utility
PRIMARY EXAMINER: Wityshyn, Michael G.
ASSISTANT EXAMINER: Gitomer, Ralph
LEGAL REPRESENTATIVE: Wegner, Cantor, Mueller & Player
NUMBER OF CLAIMS: 10
EXEMPLARY CLAIM: 1
NUMBER OF DRAWINGS: 7 Drawing Figure(s); 6 Drawing Page(s)
LINE COUNT: 1228
CAS INDEXING IS AVAILABLE FOR THIS PATENT.

AB There are described antitumour sterols, their fatty acid esters and glucosides, processes for their preparation, spontaneously dispersible agents containing these sterols, their fatty acid esters and glucosides, and their use for treating tumours.

CLM What is claimed:

3. A **pharmaceutical** composition comprising a spontaneously dispersible concentrate as claimed in claim 2, which contains 0.001 to 15% by weight of an. . . antitumor components selected from the group consisting of STIGMASTEROL-UNDECENOATE, STIGMASTEROL-DODECENOATE, STIGMASTEROL-OLEATE, STIGMASTEROL-LINOLEATE, STIGMASTEROL-LINOLENATE, .beta.-SITOSTEROL-UNDECENOATE, .beta.-SITOSTEROL-DODECENOATE, .beta.-SITOSTEROL-OLEATE, .beta.-SITOSTEROL-LINOLEATE, .beta.-SITOSTEROL-LINOLENATE, CHOLESTERYL-UNDECENOATE, CHOLESTERYL-DODECENOATE, **CHOLESTERYL-OLEATE**, CHOLESTERYL-LINOLEATE, and CHOLESTERYL-LINOLENATE.

NCL NCLM: **424/195.100**
NCLS: 536/005.000; 536/006.200; 549/408.000; 552/540.000; 552/544.000; 552/545.000; 552/547.000; 568/824.000